



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 10:34 AM UTC

PDB ID : 3OJV / pdb_00003ojv
Title : Crystal Structure of FGF1 complexed with the ectodomain of FGFR1c exhibiting an ordered ligand specificity-determining betaC'-betaE loop
Authors : Beenken, A.; Mohammadi, M.
Deposited on : 2010-08-23
Resolution : 2.60 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

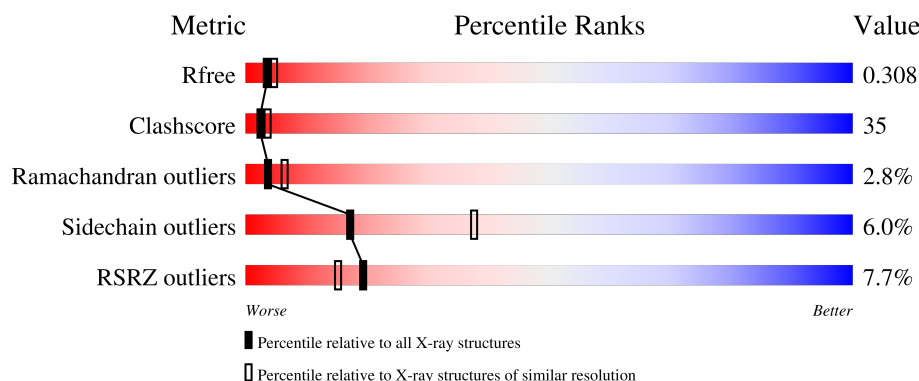
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	136	11% 46% 47% 6% 6%
1	B	136	43% 49% 6% 6%
2	C	226	11% 46% 43% 6% 6%
2	D	226	12% 47% 39% 8% 6%
3	E	6	17% 83%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	UAP	E	6	X	-	-	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5465 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Heparin-binding growth factor 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	133	Total	C	N	O	S	0	0	0
			1052	665	184	199	4			
1	B	133	Total	C	N	O	S	0	0	0
			1052	665	184	199	4			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	20	MET	-	initiating methionine	UNP P05230
B	20	MET	-	initiating methionine	UNP P05230

- Molecule 2 is a protein called Basic fibroblast growth factor receptor 1.

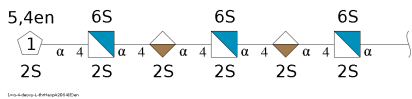
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	213	Total	C	N	O	S	0	0	0
			1608	1026	278	295	9			
2	D	213	Total	C	N	O	S	0	0	0
			1617	1028	283	297	9			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	140	MET	-	initiating methionine	UNP P11362
C	141	ALA	-	expression tag	UNP P11362
C	185	GLN	ASN	engineered mutation	UNP P11362
D	140	MET	-	initiating methionine	UNP P11362
D	141	ALA	-	expression tag	UNP P11362
D	185	GLN	ASN	engineered mutation	UNP P11362

- Molecule 3 is an oligosaccharide called 4-deoxy-2-O-sulfo-alpha-L-threo-hex-4-enopyranuronic acid-(1-4)-2-deoxy-6-O-sulfo-2-(sulfoamino)-alpha-D-glucopyranose-(1-4)-2-O-sulfo-alpha-L-idopyranuronic acid-(1-4)-2-deoxy-6-O-sulfo-2-(sulfoamino)-alpha-D-glucopyranose-(1-4)-

2-O-sulfo-alpha-L-idopyranuronic acid-(1-4)-2-deoxy-6-O-sulfo-2-(sulfoamino)-alpha-D-glucopyranose.

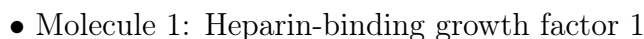


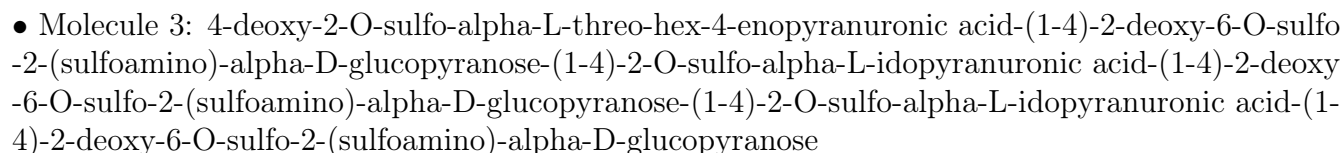
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	E	6	Total	C	N	O	S	0	0	0
			104	36	3	56	9			

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	6	Total	O	0	0
			6	6		
4	B	2	Total	O	0	0
			2	2		
4	C	13	Total	O	0	0
			13	13		
4	D	11	Total	O	0	0
			11	11		

- Molecule 1: Heparin-binding growth factor 1





Satisfaction Level	Percentage
Not at all satisfied	17%
Satisfied	83%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	53.32Å 53.42Å 80.46Å 106.44° 106.40° 94.44°	Depositor
Resolution (Å)	25.00 – 2.60 25.00 – 2.60	Depositor EDS
% Data completeness (in resolution range)	(Not available) (25.00-2.60) 95.8 (25.00-2.60)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.35 (at 2.58Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.251 , 0.309 0.253 , 0.308	Depositor DCC
R_{free} test set	2405 reflections (9.77%)	wwPDB-VP
Wilson B-factor (Å ²)	45.8	Xtriage
Anisotropy	0.124	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 48.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.018 for k,h,-h-k-l 0.478 for -k,-h,-l 0.017 for -h,-k,h+k+l	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	5465	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.95% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: SGN, IDS, UAP

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.47	0/1076	0.97	3/1453 (0.2%)
1	B	0.47	0/1076	1.00	6/1453 (0.4%)
2	C	0.50	0/1655	1.04	8/2268 (0.4%)
2	D	0.50	0/1664	1.03	9/2279 (0.4%)
All	All	0.49	0/5471	1.01	26/7453 (0.3%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
2	C	0	1

There are no bond length outliers.

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	69	VAL	N-CA-C	8.69	121.31	108.96
1	B	69	VAL	N-CA-C	8.63	121.22	108.96
2	C	279	VAL	N-CA-C	7.72	119.97	108.46
2	D	279	VAL	N-CA-C	7.54	119.70	108.46
1	B	142	GLN	N-CA-C	6.18	118.30	110.33
2	C	284	GLN	CA-C-N	-5.84	114.54	120.85
2	C	284	GLN	C-N-CA	-5.84	114.54	120.85
2	D	358	VAL	N-CA-C	5.80	116.66	108.36
2	C	358	VAL	N-CA-C	5.61	116.39	108.36
2	C	299	LYS	N-CA-C	-5.61	106.61	113.50
2	C	282	ASP	CA-C-N	-5.49	114.11	119.76
2	C	282	ASP	C-N-CA	-5.49	114.11	119.76

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	299	LYS	N-CA-C	-5.44	106.80	113.50
2	D	282	ASP	CA-C-N	-5.41	114.19	119.76
2	D	282	ASP	C-N-CA	-5.41	114.19	119.76
1	B	53	SER	N-CA-C	-5.34	106.69	114.12
1	A	72	LYS	N-CA-C	5.31	117.32	108.99
1	B	150	LEU	CA-C-N	5.28	125.26	120.03
1	B	150	LEU	C-N-CA	5.28	125.26	120.03
2	D	284	GLN	CA-C-N	-5.26	115.17	120.85
2	D	284	GLN	C-N-CA	-5.26	115.17	120.85
1	B	72	LYS	N-CA-C	5.24	117.22	108.99
1	A	53	SER	N-CA-C	-5.16	106.95	114.12
2	D	233	GLU	N-CA-C	5.05	117.65	109.06
2	C	248	VAL	N-CA-C	5.03	115.82	108.58
2	D	236	TYR	N-CA-C	5.01	118.71	112.59

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	C	280	TYR	Sidechain

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1052	0	1023	85	0
1	B	1052	0	1023	75	0
2	C	1608	0	1531	104	0
2	D	1617	0	1537	104	0
3	E	104	0	37	10	0
4	A	6	0	0	2	0
4	B	2	0	0	0	0
4	C	13	0	0	0	0
4	D	11	0	0	0	0
All	All	5465	0	5151	365	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including

hydrogen atoms). The all-atom clashscore for this structure is 35.

All (365) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:E:1:SGN:H4	3:E:2:IDS:H5	1.30	1.09
1:B:128:LYS:HD2	3:E:2:IDS:H4	1.32	1.06
1:A:123:PHE:HB2	1:A:138:THR:HG21	1.37	1.05
1:B:123:PHE:HB2	1:B:138:THR:HG21	1.37	1.02
2:D:318:THR:HG23	2:D:319:THR:H	1.33	0.93
2:C:318:THR:HG21	2:C:323:MET:HG2	1.56	0.86
2:C:265:LYS:O	2:C:356:LEU:HA	1.76	0.86
1:A:133:LYS:NZ	1:A:142:GLN:HE22	1.74	0.85
2:D:333:PHE:HA	2:D:358:VAL:HG11	1.58	0.85
2:C:333:PHE:HA	2:C:358:VAL:HG11	1.59	0.85
3:E:1:SGN:H4	3:E:2:IDS:C5	2.06	0.85
1:A:143:LYS:HD3	3:E:5:SGN:O3	1.76	0.84
2:D:265:LYS:O	2:D:356:LEU:HA	1.78	0.84
2:C:257:LEU:HD23	2:C:279:VAL:HG22	1.61	0.82
2:D:257:LEU:CD2	2:D:279:VAL:HG22	2.08	0.82
1:A:59:LEU:HD22	1:A:71:ILE:HG22	1.61	0.82
1:B:42:PRO:HD3	1:B:56:HIS:CD2	2.15	0.81
2:C:314:ALA:HB2	2:C:323:MET:HG3	1.65	0.79
2:D:183:THR:HA	2:D:184:PRO:C	2.08	0.79
2:C:291:LYS:HB2	2:C:311:LEU:HD11	1.66	0.78
2:C:257:LEU:CD2	2:C:279:VAL:HG22	2.14	0.77
2:D:277:CYS:HG	2:D:341:CYS:HG	0.91	0.77
2:D:222:PRO:O	2:D:225:LYS:HG2	1.85	0.77
2:D:209:ARG:HG2	2:D:209:ARG:HH11	1.49	0.75
2:D:277:CYS:O	2:D:324:GLU:HG3	1.86	0.75
2:D:314:ALA:HB2	2:D:323:MET:HG3	1.69	0.74
1:A:134:ARG:C	1:A:136:PRO:HD2	2.14	0.73
2:C:183:THR:HA	2:C:184:PRO:C	2.13	0.73
1:A:133:LYS:HZ3	1:A:142:GLN:NE2	1.88	0.72
2:D:257:LEU:HD23	2:D:279:VAL:HG22	1.72	0.72
2:D:336:ALA:HB2	2:D:358:VAL:HG23	1.73	0.71
2:C:292:HIS:HD2	2:C:338:GLU:HB3	1.56	0.70
2:D:318:THR:HG23	2:D:319:THR:N	2.07	0.70
3:E:1:SGN:C4	3:E:2:IDS:H5	2.14	0.69
1:B:119:GLU:C	1:B:121:ASN:H	1.99	0.69
1:B:59:LEU:HD22	1:B:71:ILE:HG22	1.74	0.69
2:C:262:PRO:HD2	2:C:352:HIS:HB3	1.74	0.68
1:B:139:HIS:HB3	1:B:142:GLN:HG3	1.73	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:187:THR:HG22	2:D:188:LEU:N	2.08	0.68
1:A:119:GLU:C	1:A:121:ASN:H	2.02	0.68
2:C:273:VAL:HG13	2:C:328:LEU:HD12	1.77	0.67
2:D:291:LYS:HB2	2:D:311:LEU:HD11	1.76	0.67
2:C:336:ALA:HB2	2:C:358:VAL:HG23	1.76	0.67
1:B:133:LYS:NZ	1:B:142:GLN:HE22	1.93	0.67
1:A:26:PRO:O	1:A:27:LYS:HD2	1.95	0.67
2:D:292:HIS:HD2	2:D:308:VAL:HG12	1.59	0.67
1:A:76:THR:OG1	1:A:78:GLN:HG3	1.96	0.66
1:B:76:THR:OG1	1:B:78:GLN:HG3	1.96	0.65
2:C:188:LEU:HD11	2:C:190:TRP:NE1	2.11	0.65
2:D:261:LEU:CD1	2:D:278:LYS:HG3	2.27	0.65
2:D:269:LEU:H	2:D:269:LEU:HD23	1.61	0.65
2:D:294:GLU:HG2	2:D:299:LYS:HG2	1.76	0.65
1:A:133:LYS:NZ	1:A:142:GLN:NE2	2.44	0.65
2:C:282:ASP:HB3	2:C:283:PRO:CD	2.25	0.65
1:A:71:ILE:HD12	1:A:80:LEU:HD23	1.79	0.65
1:A:133:LYS:HZ3	1:A:142:GLN:CD	2.05	0.65
1:A:22:ASN:O	1:A:27:LYS:NZ	2.30	0.65
2:C:292:HIS:CD2	2:C:338:GLU:HB3	2.32	0.64
1:B:133:LYS:NZ	1:B:142:GLN:NE2	2.45	0.64
2:C:187:THR:HG22	2:C:188:LEU:N	2.10	0.64
1:A:33:ASN:ND2	1:A:128:LYS:HG2	2.13	0.63
1:B:127:LYS:HD2	1:B:131:SER:OG	1.97	0.63
2:D:316:VAL:O	2:D:317:ASN:CB	2.45	0.63
1:B:54:ASP:OD1	1:B:56:HIS:N	2.27	0.63
1:A:29:LEU:HD12	1:A:59:LEU:HD13	1.79	0.63
1:A:42:PRO:HD3	1:A:56:HIS:CD2	2.33	0.63
2:D:317:ASN:O	2:D:319:THR:N	2.32	0.62
1:B:71:ILE:HD12	1:B:80:LEU:HD23	1.81	0.62
2:D:269:LEU:HD23	2:D:269:LEU:N	2.15	0.62
2:D:188:LEU:HD23	2:D:189:ARG:N	2.13	0.62
1:B:26:PRO:O	1:B:27:LYS:HD2	1.99	0.62
1:A:34:GLY:HA3	1:A:36:HIS:CE1	2.34	0.62
1:B:134:ARG:C	1:B:136:PRO:HD2	2.24	0.62
2:D:287:ILE:N	2:D:287:ILE:HD12	2.14	0.62
2:D:290:LEU:HD22	2:D:310:ILE:HD13	1.81	0.62
2:C:312:LYS:HG2	2:C:323:MET:HE1	1.82	0.62
2:C:316:VAL:O	2:C:317:ASN:CB	2.48	0.61
2:C:317:ASN:O	2:C:319:THR:N	2.32	0.61
2:D:257:LEU:HD21	2:D:279:VAL:HG22	1.81	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:282:ASP:HB3	2:C:283:PRO:HD2	1.81	0.61
1:A:133:LYS:HZ1	1:A:142:GLN:HE22	1.46	0.61
2:C:159:GLU:C	2:C:161:MET:H	2.09	0.61
1:A:102:GLU:HG3	1:A:112:TYR:CE1	2.35	0.61
2:C:187:THR:HG22	2:C:188:LEU:H	1.66	0.61
1:A:50:ARG:HE	2:C:163:LYS:HA	1.65	0.61
2:C:147:ASN:N	2:C:147:ASN:HD22	1.97	0.60
1:B:22:ASN:O	1:B:27:LYS:NZ	2.34	0.60
1:B:133:LYS:CE	1:B:142:GLN:HE22	2.14	0.60
2:D:209:ARG:HG2	2:D:209:ARG:NH1	2.12	0.60
2:D:262:PRO:HD2	2:D:352:HIS:HB3	1.82	0.60
2:D:261:LEU:HD12	2:D:278:LYS:HG3	1.83	0.60
2:D:159:GLU:C	2:D:161:MET:H	2.10	0.60
2:D:295:VAL:HG22	2:D:296:ASN:ND2	2.17	0.60
1:A:54:ASP:OD1	1:A:55:GLN:N	2.33	0.60
2:D:318:THR:CG2	2:D:319:THR:H	2.11	0.60
1:B:123:PHE:CB	1:B:138:THR:HG21	2.24	0.60
2:D:261:LEU:HB3	2:D:276:MET:O	2.02	0.59
2:D:282:ASP:HB3	2:D:283:PRO:CD	2.32	0.59
1:A:127:LYS:HD2	1:A:131:SER:OG	2.02	0.59
2:D:163:LYS:HD2	2:D:166:HIS:CE1	2.36	0.59
1:B:82:MET:HA	1:B:87:LEU:O	2.03	0.59
1:A:135:GLY:N	1:A:136:PRO:HD2	2.18	0.59
1:A:40:ILE:O	1:A:56:HIS:HB3	2.02	0.58
2:C:225:LYS:HE2	2:C:246:ASP:OD1	2.02	0.58
2:C:318:THR:O	2:C:319:THR:HG23	2.03	0.58
1:A:119:GLU:C	1:A:121:ASN:N	2.61	0.58
2:C:155:TRP:HB3	2:C:161:MET:HE1	1.85	0.58
1:A:33:ASN:HD21	1:A:128:LYS:HG2	1.69	0.58
1:A:70:TYR:CE2	2:C:316:VAL:HB	2.39	0.58
1:A:82:MET:HA	1:A:87:LEU:O	2.04	0.58
2:D:318:THR:O	2:D:319:THR:HG23	2.04	0.58
2:D:344:GLY:HA3	2:D:349:LEU:HD23	1.86	0.58
1:A:123:PHE:HD1	1:A:138:THR:HG22	1.68	0.57
1:A:39:ARG:NE	1:A:41:LEU:HD21	2.18	0.57
2:C:287:ILE:HD12	2:C:287:ILE:N	2.19	0.57
1:A:95:ASN:N	1:A:95:ASN:ND2	2.53	0.57
1:A:120:LYS:HB3	1:A:122:TRP:NE1	2.20	0.57
1:B:40:ILE:O	1:B:56:HIS:HB3	2.04	0.57
1:B:119:GLU:C	1:B:121:ASN:N	2.59	0.57
1:B:122:TRP:HB3	1:B:135:GLY:HA3	1.86	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:261:LEU:HB3	2:C:276:MET:O	2.05	0.56
2:D:190:TRP:CZ3	2:D:230:CYS:HB3	2.40	0.56
2:D:264:ASN:O	2:D:265:LYS:HG3	2.05	0.56
2:D:292:HIS:ND1	2:D:338:GLU:HB3	2.20	0.56
2:D:316:VAL:O	2:D:317:ASN:HB2	2.06	0.56
2:D:277:CYS:HB2	2:D:289:TRP:CH2	2.41	0.56
2:D:317:ASN:OD1	2:D:318:THR:N	2.38	0.55
2:D:187:THR:CG2	2:D:188:LEU:N	2.68	0.55
1:A:84:THR:HA	1:A:117:HIS:NE2	2.21	0.55
1:A:95:ASN:N	1:A:95:ASN:HD22	2.03	0.55
1:B:39:ARG:O	1:B:46:VAL:HG13	2.07	0.55
2:C:174:VAL:HG23	2:C:220:VAL:HG11	1.88	0.55
1:A:133:LYS:O	1:A:134:ARG:C	2.50	0.54
2:C:301:GLY:HA2	2:C:307:TYR:CE1	2.42	0.54
2:C:252:PRO:C	2:C:347:ILE:HD13	2.33	0.54
1:A:128:LYS:HG3	3:E:4:IDS:O1S	2.07	0.54
1:A:82:MET:HE3	1:A:88:LEU:HD21	1.90	0.54
2:D:264:ASN:HD22	2:D:264:ASN:H	1.56	0.53
2:D:277:CYS:HG	2:D:341:CYS:CB	2.22	0.53
2:D:333:PHE:HA	2:D:358:VAL:CG1	2.35	0.53
1:A:133:LYS:O	1:A:134:ARG:O	2.25	0.53
1:B:135:GLY:N	1:B:136:PRO:HD2	2.24	0.53
1:B:120:LYS:HB3	1:B:122:TRP:NE1	2.23	0.53
1:B:133:LYS:O	1:B:134:ARG:C	2.51	0.53
1:B:133:LYS:O	1:B:134:ARG:O	2.27	0.53
2:D:314:ALA:CB	2:D:323:MET:HG3	2.39	0.53
2:C:265:LYS:NZ	2:C:274:GLU:O	2.40	0.53
1:B:133:LYS:O	1:B:133:LYS:HG2	2.08	0.53
2:C:264:ASN:O	2:C:265:LYS:HG3	2.10	0.52
1:A:59:LEU:HD22	1:A:71:ILE:CG2	2.38	0.52
2:C:277:CYS:HG	2:C:341:CYS:CB	2.21	0.52
2:C:295:VAL:O	2:C:296:ASN:HB2	2.08	0.52
2:C:317:ASN:OD1	2:C:318:THR:N	2.37	0.52
2:C:333:PHE:HA	2:C:358:VAL:CG1	2.36	0.52
2:D:147:ASN:N	2:D:147:ASN:HD22	2.06	0.52
2:D:159:GLU:O	2:D:161:MET:N	2.42	0.52
2:C:159:GLU:O	2:C:161:MET:N	2.42	0.52
1:B:133:LYS:HE2	1:B:142:GLN:HE22	1.75	0.52
2:C:314:ALA:CB	2:C:323:MET:HG3	2.38	0.52
2:D:293:ILE:CG2	2:D:294:GLU:N	2.71	0.52
2:D:294:GLU:HG2	2:D:299:LYS:CG	2.40	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:273:VAL:CG1	2:C:328:LEU:HD12	2.40	0.52
2:C:317:ASN:CG	2:C:318:THR:H	2.16	0.52
1:B:97:GLU:HG2	1:B:116:LYS:CD	2.41	0.51
2:D:273:VAL:CG1	2:D:328:LEU:HD12	2.40	0.51
2:C:172:LYS:O	2:C:220:VAL:HG13	2.10	0.51
2:D:161:MET:HE3	2:D:178:CYS:HA	1.91	0.51
2:D:245:LEU:HD22	2:D:246:ASP:N	2.26	0.51
2:C:358:VAL:O	2:C:359:LEU:HG	2.10	0.51
2:D:187:THR:HG22	2:D:188:LEU:H	1.75	0.51
2:D:273:VAL:HG13	2:D:328:LEU:HD12	1.92	0.51
2:C:273:VAL:HG12	2:C:331:VAL:HG21	1.91	0.51
1:B:89:TYR:CE1	1:B:91:SER:HB2	2.46	0.51
2:C:257:LEU:HD21	2:C:279:VAL:HG22	1.93	0.51
2:C:170:ALA:O	2:C:171:ALA:HB3	2.10	0.51
1:A:135:GLY:O	1:A:138:THR:HB	2.11	0.51
2:C:186:PRO:HG2	2:C:213:TRP:HH2	1.75	0.51
2:C:158:PRO:O	2:C:161:MET:HB2	2.11	0.51
1:A:128:LYS:HZ1	3:E:5:SGN:H3	1.76	0.50
2:C:186:PRO:HG2	2:C:213:TRP:CH2	2.46	0.50
1:A:122:TRP:HB3	1:A:135:GLY:HA3	1.92	0.50
2:C:290:LEU:HD22	2:C:310:ILE:CD1	2.40	0.50
2:D:312:LYS:HG2	2:D:323:MET:HE1	1.92	0.50
2:D:267:VAL:HG22	2:D:356:LEU:HD11	1.93	0.50
1:B:70:TYR:CE2	2:D:316:VAL:HB	2.46	0.50
1:B:89:TYR:HE1	1:B:91:SER:HB2	1.76	0.50
1:B:148:LEU:O	1:B:150:LEU:HD23	2.11	0.50
2:C:303:ASP:OD2	2:C:305:LEU:HD12	2.12	0.50
1:B:133:LYS:HZ3	1:B:142:GLN:NE2	2.09	0.50
1:B:135:GLY:O	1:B:138:THR:HB	2.11	0.49
1:A:123:PHE:CB	1:A:138:THR:HG21	2.26	0.49
2:D:187:THR:CG2	2:D:188:LEU:H	2.24	0.49
1:B:83:ASP:OD2	1:B:83:ASP:C	2.55	0.49
2:C:312:LYS:HE2	2:C:323:MET:HE2	1.95	0.49
1:B:36:HIS:NE2	1:B:130:GLY:HA3	2.28	0.49
2:C:277:CYS:HB2	2:C:289:TRP:CH2	2.48	0.49
2:D:255:PRO:O	2:D:350:SER:OG	2.24	0.48
2:C:187:THR:CG2	2:C:188:LEU:N	2.77	0.48
2:C:292:HIS:HD2	2:C:338:GLU:CB	2.26	0.48
1:B:24:LYS:HG2	2:D:320:ASP:HB3	1.95	0.48
2:C:290:LEU:HD22	2:C:310:ILE:HD13	1.95	0.48
2:C:342:LEU:HD21	2:C:349:LEU:HB2	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:40:ILE:HA	1:B:46:VAL:HG22	1.95	0.48
1:B:52:ARG:O	1:B:52:ARG:HG2	2.14	0.48
1:B:30:TYR:HB3	1:B:150:LEU:HD23	1.96	0.48
1:B:54:ASP:OD1	1:B:55:GLN:N	2.47	0.48
1:A:39:ARG:O	1:A:46:VAL:HG13	2.13	0.48
2:D:282:ASP:HB3	2:D:283:PRO:HD2	1.96	0.48
1:A:21:GLY:N	1:A:27:LYS:HE3	2.29	0.47
2:C:270:GLY:HA2	2:C:330:ASN:ND2	2.29	0.47
2:C:293:ILE:CG2	2:C:294:GLU:N	2.75	0.47
2:C:302:PRO:O	2:C:303:ASP:HB3	2.12	0.47
2:D:344:GLY:HA3	2:D:349:LEU:CD2	2.44	0.47
1:A:127:LYS:HB2	1:A:129:ASN:OD1	2.14	0.47
2:C:220:VAL:HG23	2:C:247:VAL:HG13	1.95	0.47
2:D:170:ALA:O	2:D:171:ALA:HB3	2.14	0.47
1:A:36:HIS:HE1	1:A:128:LYS:O	1.97	0.47
1:A:95:ASN:O	1:A:98:CYS:HB2	2.14	0.47
1:B:84:THR:HA	1:B:117:HIS:NE2	2.29	0.47
1:A:123:PHE:HB2	1:A:138:THR:CG2	2.27	0.47
1:A:33:ASN:HD22	1:A:128:LYS:HA	1.78	0.47
1:A:52:ARG:O	1:A:52:ARG:HG2	2.14	0.47
1:A:139:HIS:HB3	1:A:142:GLN:HG3	1.97	0.47
2:C:151:VAL:HB	2:C:183:THR:H	1.79	0.47
2:C:190:TRP:CZ3	2:C:230:CYS:HB3	2.49	0.47
2:C:275:PHE:O	2:C:325:VAL:HA	2.15	0.47
2:D:275:PHE:O	2:D:325:VAL:HA	2.15	0.47
1:B:51:ASP:C	1:B:53:SER:H	2.23	0.47
1:B:50:ARG:HE	2:D:163:LYS:HA	1.80	0.47
1:B:50:ARG:NE	2:D:163:LYS:HA	2.30	0.47
2:D:265:LYS:O	2:D:356:LEU:HD12	2.15	0.47
1:A:103:ARG:HG3	4:A:4015:HOH:O	2.15	0.46
1:A:26:PRO:C	1:A:27:LYS:HD2	2.39	0.46
2:C:286:HIS:C	2:C:287:ILE:HD12	2.40	0.46
2:D:264:ASN:H	2:D:264:ASN:ND2	2.14	0.46
2:D:151:VAL:HB	2:D:183:THR:H	1.80	0.46
1:A:139:HIS:H	1:A:142:GLN:HG3	1.80	0.46
2:C:159:GLU:C	2:C:161:MET:N	2.74	0.46
2:C:183:THR:CA	2:C:184:PRO:C	2.88	0.46
2:C:153:PRO:HA	2:C:181:SER:O	2.15	0.46
2:D:159:GLU:C	2:D:161:MET:N	2.74	0.46
2:D:271:SER:O	2:D:331:VAL:HG23	2.16	0.46
1:A:22:ASN:O	1:A:24:LYS:N	2.49	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:102:GLU:HG3	1:B:112:TYR:CE1	2.50	0.46
1:A:128:LYS:NZ	3:E:4:IDS:O6A	2.48	0.45
1:A:139:HIS:O	1:A:142:GLN:HB2	2.16	0.45
1:B:84:THR:O	1:B:84:THR:HG22	2.16	0.45
1:B:76:THR:CB	1:B:78:GLN:HG3	2.47	0.45
2:C:187:THR:CG2	2:C:188:LEU:H	2.29	0.45
1:B:113:ILE:HG23	1:B:122:TRP:O	2.16	0.45
2:D:153:PRO:HA	2:D:181:SER:O	2.16	0.45
1:A:28:LEU:HD11	1:A:152:VAL:HA	1.99	0.45
1:B:97:GLU:HG2	1:B:116:LYS:HD2	1.98	0.45
1:A:51:ASP:C	1:A:53:SER:H	2.25	0.45
1:B:119:GLU:HA	1:B:119:GLU:OE1	2.16	0.45
1:B:120:LYS:O	1:B:122:TRP:HD1	2.00	0.45
2:D:168:VAL:HG22	2:D:169:PRO:HD2	1.98	0.45
2:D:290:LEU:CD2	2:D:310:ILE:HD13	2.46	0.45
2:C:245:LEU:HD22	2:C:246:ASP:N	2.32	0.45
2:C:295:VAL:O	2:C:295:VAL:HG13	2.16	0.45
2:D:206:TYR:HE1	2:D:208:VAL:CG2	2.29	0.45
1:B:101:LEU:HD21	1:B:115:LYS:HG2	1.98	0.45
2:C:155:TRP:CE2	2:C:239:ILE:HD12	2.52	0.45
2:C:334:GLU:C	2:C:336:ALA:H	2.25	0.45
1:A:64:GLU:OE1	1:A:64:GLU:O	2.35	0.45
2:C:176:PHE:HB2	2:C:215:ILE:HG22	1.99	0.45
2:C:356:LEU:C	2:C:356:LEU:HD23	2.41	0.45
2:D:220:VAL:HG23	2:D:247:VAL:HG13	1.98	0.45
2:D:315:GLY:O	2:D:318:THR:HG22	2.16	0.45
3:E:3:SGN:H62	3:E:4:IDS:C1	2.47	0.45
1:B:21:GLY:N	1:B:27:LYS:HE3	2.32	0.44
2:C:312:LYS:HE2	2:C:323:MET:CE	2.47	0.44
1:A:81:ALA:O	1:A:88:LEU:HA	2.17	0.44
2:C:251:SER:HB2	2:C:282:ASP:HB2	1.98	0.44
1:B:81:ALA:O	1:B:88:LEU:HA	2.18	0.44
2:C:234:ASN:OD1	2:C:234:ASN:C	2.60	0.44
2:D:312:LYS:HE2	2:D:323:MET:HE2	1.99	0.44
2:D:344:GLY:CA	2:D:349:LEU:HD23	2.47	0.44
1:A:33:ASN:C	1:A:35:GLY:H	2.26	0.44
1:A:138:THR:HG23	1:A:145:ILE:CG2	2.48	0.44
2:D:318:THR:O	2:D:319:THR:OG1	2.33	0.44
1:A:39:ARG:O	1:A:46:VAL:HA	2.18	0.44
1:A:39:ARG:NH2	1:A:47:ASP:OD2	2.50	0.44
1:B:37:PHE:CE2	1:B:52:ARG:HB2	2.53	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:39:ARG:O	1:B:46:VAL:HA	2.17	0.44
2:D:315:GLY:O	2:D:316:VAL:C	2.60	0.44
1:A:76:THR:CB	1:A:78:GLN:HG3	2.47	0.44
1:A:133:LYS:O	1:A:133:LYS:HG2	2.18	0.44
2:C:188:LEU:HD23	2:C:232:VAL:HG22	2.00	0.44
2:D:302:PRO:O	2:D:303:ASP:HB3	2.18	0.44
1:B:42:PRO:C	1:B:44:GLY:H	2.26	0.44
2:D:174:VAL:HG23	2:D:220:VAL:HG11	2.00	0.44
2:C:318:THR:HG21	2:C:323:MET:CE	2.48	0.44
1:A:84:THR:O	1:A:122:TRP:HH2	2.00	0.43
1:A:120:LYS:O	1:A:122:TRP:HD1	2.01	0.43
1:B:42:PRO:HD3	1:B:56:HIS:HD2	1.76	0.43
2:D:334:GLU:C	2:D:336:ALA:H	2.26	0.43
1:B:26:PRO:C	1:B:27:LYS:HD2	2.43	0.43
2:C:188:LEU:HD11	2:C:190:TRP:HE1	1.80	0.43
2:D:295:VAL:O	2:D:296:ASN:HB2	2.18	0.43
1:B:22:ASN:O	1:B:24:LYS:N	2.51	0.43
1:B:39:ARG:NE	1:B:41:LEU:HD21	2.33	0.43
1:B:86:GLY:O	1:B:134:ARG:HA	2.18	0.43
2:D:190:TRP:CH2	2:D:230:CYS:HB3	2.53	0.43
2:D:288:GLN:HE21	2:D:288:GLN:HB3	1.48	0.43
1:A:42:PRO:C	1:A:44:GLY:H	2.26	0.43
1:B:33:ASN:C	1:B:35:GLY:H	2.25	0.43
2:C:289:TRP:CD2	2:C:326:LEU:HB2	2.53	0.43
2:D:270:GLY:HA2	2:D:330:ASN:ND2	2.32	0.43
1:A:61:LEU:N	1:A:61:LEU:HD12	2.34	0.43
1:A:24:LYS:HE2	4:A:4002:HOH:O	2.18	0.43
2:D:295:VAL:O	2:D:295:VAL:HG13	2.17	0.43
1:A:86:GLY:O	1:A:134:ARG:HA	2.19	0.43
1:A:61:LEU:HB3	2:C:284:GLN:HG3	2.01	0.43
2:D:186:PRO:HG2	2:D:213:TRP:CH2	2.54	0.43
1:A:39:ARG:HH21	1:A:47:ASP:CG	2.27	0.42
2:C:330:ASN:O	2:C:331:VAL:C	2.62	0.42
1:B:119:GLU:O	1:B:121:ASN:N	2.52	0.42
1:A:37:PHE:CE2	1:A:52:ARG:HB2	2.54	0.42
2:D:172:LYS:O	2:D:220:VAL:HG13	2.19	0.42
2:D:330:ASN:O	2:D:331:VAL:C	2.62	0.42
1:B:110:ASN:HB3	1:B:112:TYR:CE2	2.54	0.42
1:A:83:ASP:C	1:A:83:ASP:OD2	2.61	0.42
1:B:113:ILE:HD13	1:B:121:ASN:ND2	2.33	0.42
2:C:255:PRO:O	2:C:350:SER:OG	2.35	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:264:ASN:C	2:C:265:LYS:HG3	2.45	0.42
2:C:252:PRO:HA	2:C:347:ILE:CD1	2.50	0.42
2:D:358:VAL:O	2:D:359:LEU:HD23	2.20	0.42
1:B:95:ASN:H	1:B:98:CYS:HB2	1.85	0.42
1:B:123:PHE:HD1	1:B:138:THR:HG22	1.85	0.42
2:C:186:PRO:HB2	2:C:232:VAL:HG13	2.02	0.42
1:A:95:ASN:H	1:A:98:CYS:HB2	1.84	0.42
1:B:41:LEU:O	1:B:44:GLY:N	2.49	0.42
2:C:231:ILE:HG22	2:C:233:GLU:HG2	2.00	0.42
1:B:24:LYS:O	1:B:60:GLN:NE2	2.52	0.42
2:C:282:ASP:CB	2:C:283:PRO:CD	2.96	0.42
2:C:316:VAL:O	2:C:317:ASN:HB2	2.19	0.42
2:D:183:THR:CA	2:D:184:PRO:C	2.84	0.42
2:D:323:MET:C	2:D:325:VAL:H	2.28	0.42
2:C:286:HIS:CD2	2:C:314:ALA:HB3	2.55	0.41
2:C:182:GLY:O	2:C:185:GLN:HA	2.19	0.41
2:D:333:PHE:CA	2:D:358:VAL:HG11	2.41	0.41
2:C:316:VAL:O	2:C:317:ASN:HB3	2.18	0.41
1:B:30:TYR:CD1	1:B:30:TYR:C	2.98	0.41
2:C:323:MET:C	2:C:325:VAL:H	2.28	0.41
2:C:206:TYR:HE1	2:C:208:VAL:CG2	2.34	0.41
2:D:305:LEU:HA	2:D:306:PRO:HD2	1.89	0.41
2:D:191:LEU:HD23	2:D:196:GLU:HA	2.01	0.41
2:C:147:ASN:N	2:C:147:ASN:ND2	2.68	0.41
1:A:54:ASP:OD1	1:A:56:HIS:N	2.42	0.41
1:A:41:LEU:O	1:A:44:GLY:N	2.49	0.41
2:C:188:LEU:HD21	2:C:230:CYS:HB2	2.03	0.41
2:C:318:THR:HG21	2:C:323:MET:HE2	2.02	0.41
2:D:286:HIS:C	2:D:287:ILE:HD12	2.46	0.41
2:D:318:THR:O	2:D:319:THR:CB	2.69	0.41
1:A:123:PHE:CD1	1:A:138:THR:HG22	2.50	0.40
1:B:84:THR:O	1:B:122:TRP:HH2	2.04	0.40
1:B:72:LYS:HD2	1:B:79:TYR:CZ	2.56	0.40
2:C:315:GLY:O	2:C:316:VAL:C	2.60	0.40
1:A:82:MET:HE3	1:A:88:LEU:CD2	2.51	0.40
2:C:254:ARG:H	2:C:254:ARG:HG2	1.60	0.40
2:C:323:MET:C	2:C:325:VAL:N	2.79	0.40
2:D:282:ASP:CB	2:D:283:PRO:CD	2.99	0.40
1:A:41:LEU:HD12	1:A:45:THR:HB	2.04	0.40
1:A:143:LYS:HD3	3:E:5:SGN:HO3	1.82	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	131/136 (96%)	113 (86%)	15 (12%)	3 (2%)	5	9
1	B	131/136 (96%)	113 (86%)	14 (11%)	4 (3%)	3	5
2	C	211/226 (93%)	184 (87%)	21 (10%)	6 (3%)	4	6
2	D	211/226 (93%)	185 (88%)	20 (10%)	6 (3%)	4	6
All	All	684/724 (94%)	595 (87%)	70 (10%)	19 (3%)	4	6

All (19) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	33	ASN
1	A	134	ARG
1	B	33	ASN
1	B	134	ARG
2	C	160	LYS
2	C	303	ASP
2	C	317	ASN
2	C	318	THR
2	C	319	THR
2	D	160	LYS
2	D	303	ASP
2	D	317	ASN
2	D	318	THR
2	D	319	THR
1	A	23	TYR
1	B	23	TYR
2	C	296	ASN
2	D	296	ASN
1	B	52	ARG

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	113/119 (95%)	110 (97%)	3 (3%)	39	67
1	B	113/119 (95%)	107 (95%)	6 (5%)	20	43
2	C	170/198 (86%)	159 (94%)	11 (6%)	15	35
2	D	171/198 (86%)	157 (92%)	14 (8%)	10	24
All	All	567/634 (89%)	533 (94%)	34 (6%)	17	37

All (34) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	64	GLU
1	A	95	ASN
1	A	121	ASN
1	B	33	ASN
1	B	66	VAL
1	B	69	VAL
1	B	106	GLU
1	B	121	ASN
1	B	152	VAL
2	C	147	ASN
2	C	156	THR
2	C	168	VAL
2	C	173	THR
2	C	180	SER
2	C	233	GLU
2	C	245	LEU
2	C	320	ASP
2	C	342	LEU
2	C	351	HIS
2	C	355	TRP
2	D	168	VAL
2	D	188	LEU
2	D	245	LEU
2	D	269	LEU

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Mol	Chain	Res	Type
2	D	288	GLN
2	D	293	ILE
2	D	294	GLU
2	D	304	ASN
2	D	313	THR
2	D	333	PHE
2	D	334	GLU
2	D	342	LEU
2	D	347	ILE
2	D	355	TRP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	33	ASN
1	A	60	GLN
1	A	78	GLN
1	A	95	ASN
1	A	121	ASN
1	A	142	GLN
1	B	60	GLN
1	B	121	ASN
1	B	142	GLN
2	C	185	GLN
2	C	227	ASN
2	C	284	GLN
2	C	288	GLN
2	C	330	ASN
2	D	166	HIS
2	D	185	GLN
2	D	201	HIS
2	D	227	ASN
2	D	264	ASN
2	D	288	GLN
2	D	296	ASN
2	D	330	ASN
2	D	351	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

6 monosaccharides are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	SGN	E	1	3	19,19,20	2.37	5 (26%)	23,29,31	1.61	5 (21%)
3	IDS	E	2	3	16,16,17	1.91	6 (37%)	16,24,26	1.40	1 (6%)
3	SGN	E	3	3	19,19,20	2.10	4 (21%)	23,29,31	2.01	7 (30%)
3	IDS	E	4	3	16,16,17	2.00	4 (25%)	16,24,26	2.71	4 (25%)
3	SGN	E	5	3	19,19,20	2.43	4 (21%)	23,29,31	1.84	7 (30%)
3	UAP	E	6	3	15,15,16	2.95	8 (53%)	18,22,24	1.48	3 (16%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	SGN	E	1	3	-	2/11/28/31	0/1/1/1
3	IDS	E	2	3	-	0/9/26/29	0/1/1/1
3	SGN	E	3	3	-	1/11/28/31	0/1/1/1
3	IDS	E	4	3	-	0/9/26/29	0/1/1/1
3	SGN	E	5	3	-	0/11/28/31	0/1/1/1
3	UAP	E	6	3	1/1/5/7	0/9/22/25	0/1/1/1

All (31) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	E	5	SGN	S1-N2	8.42	1.70	1.59
3	E	6	UAP	O6A-C6	6.21	1.37	1.22

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	E	1	SGN	O2S-S1	6.04	1.49	1.42
3	E	3	SGN	O2S-S1	5.68	1.48	1.42
3	E	6	UAP	C1-C2	5.17	1.60	1.51
3	E	4	IDS	C3-C2	4.83	1.63	1.53
3	E	1	SGN	O1S-S1	4.80	1.47	1.42
3	E	3	SGN	S1-N2	4.61	1.65	1.59
3	E	6	UAP	C3-C4	4.35	1.55	1.50
3	E	6	UAP	C2-C3	3.66	1.59	1.52
3	E	1	SGN	C2-N2	-3.59	1.41	1.47
3	E	4	IDS	O5-C1	3.53	1.49	1.43
3	E	2	IDS	C1-C2	3.52	1.57	1.51
3	E	1	SGN	S1-N2	3.52	1.63	1.59
3	E	2	IDS	C4-C3	3.37	1.61	1.52
3	E	5	SGN	O5-C1	3.36	1.49	1.43
3	E	3	SGN	C2-N2	-3.08	1.42	1.47
3	E	1	SGN	O4-C4	2.96	1.50	1.43
3	E	6	UAP	O2-C2	2.92	1.52	1.47
3	E	2	IDS	C3-C2	2.90	1.59	1.53
3	E	4	IDS	O5-C5	2.70	1.48	1.43
3	E	2	IDS	O5-C1	2.67	1.48	1.43
3	E	5	SGN	O5-C5	2.66	1.48	1.43
3	E	2	IDS	C4-C5	2.65	1.57	1.53
3	E	6	UAP	O3S-S	2.57	1.56	1.45
3	E	2	IDS	O5-C5	2.36	1.47	1.43
3	E	6	UAP	O6B-C6	-2.34	1.24	1.30
3	E	3	SGN	O5-C1	2.23	1.47	1.43
3	E	5	SGN	C4-C3	2.11	1.57	1.52
3	E	4	IDS	C4-C3	2.06	1.57	1.52
3	E	6	UAP	C4-C5	2.04	1.36	1.33

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	E	4	IDS	O2-C2-C3	8.87	119.34	106.95
3	E	3	SGN	C4-C3-C2	-5.55	102.88	111.02
3	E	3	SGN	O6-C6-C5	3.82	114.38	107.57
3	E	2	IDS	O2-C2-C3	3.76	112.20	106.95
3	E	5	SGN	O6-C6-C5	3.64	114.06	107.57
3	E	1	SGN	O2S-S1-N2	-3.61	102.84	108.88
3	E	6	UAP	O6B-C6-C5	3.38	121.72	114.06
3	E	5	SGN	C3-C2-N2	3.29	114.64	110.32
3	E	6	UAP	O1S-S-O2S	3.27	120.00	108.56

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	E	3	SGN	O5-C1-C2	-3.19	106.36	111.29
3	E	5	SGN	C4-C3-C2	-3.17	106.38	111.02
3	E	5	SGN	O1S-S1-N2	-3.14	103.63	108.88
3	E	1	SGN	O6-C6-C5	-2.95	102.31	107.57
3	E	3	SGN	C1-C2-N2	-2.84	105.73	110.22
3	E	1	SGN	C4-C3-C2	-2.76	106.98	111.02
3	E	5	SGN	O5S-S2-O4S	2.68	122.59	112.24
3	E	1	SGN	O5S-S2-O4S	2.64	122.41	112.24
3	E	5	SGN	C1-O5-C5	2.60	115.67	112.19
3	E	5	SGN	C1-C2-N2	-2.59	106.13	110.22
3	E	4	IDS	O2S-S-O1S	2.57	122.16	112.24
3	E	3	SGN	O5S-S2-O4S	2.41	121.54	112.24
3	E	4	IDS	O3-C3-C4	-2.37	104.80	110.38
3	E	4	IDS	O6B-C6-O6A	-2.31	118.85	124.08
3	E	3	SGN	C3-C2-N2	2.24	113.26	110.32
3	E	3	SGN	O2S-S1-N2	-2.14	105.30	108.88
3	E	1	SGN	O1S-S1-N2	-2.05	105.45	108.88
3	E	6	UAP	O6B-C6-O6A	-2.01	119.11	123.90

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	E	6	UAP	C1

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	E	1	SGN	C6-O6-S2-O6S
3	E	1	SGN	C6-O6-S2-O4S
3	E	3	SGN	C6-O6-S2-O5S

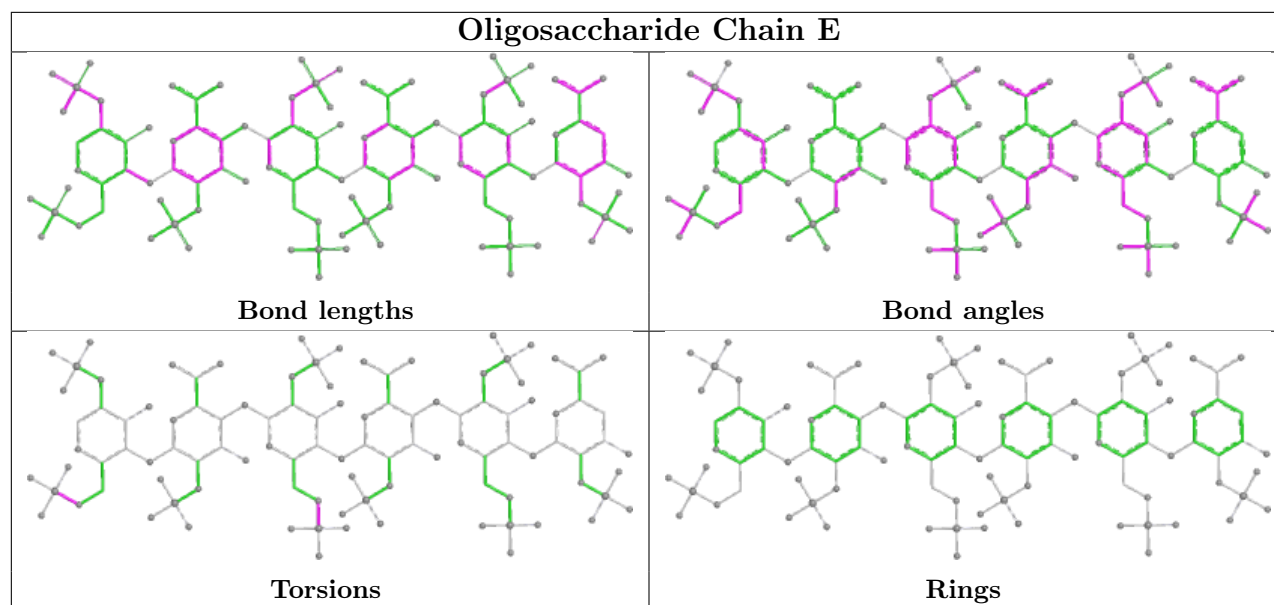
There are no ring outliers.

5 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	E	1	SGN	3	0
3	E	2	IDS	4	0
3	E	3	SGN	1	0
3	E	5	SGN	3	0
3	E	4	IDS	3	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths,

bond angles, torsion angles, and ring geometry for oligosaccharide.



5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	133/136 (97%)	0.51	1 (0%) 82 80	21, 44, 60, 63	0
1	B	133/136 (97%)	0.58	2 (1%) 72 68	23, 44, 59, 65	0
2	C	213/226 (94%)	0.52	24 (11%) 10 8	19, 38, 74, 79	0
2	D	213/226 (94%)	0.53	26 (12%) 8 6	20, 37, 74, 79	0
All	All	692/724 (95%)	0.53	53 (7%) 19 15	19, 40, 69, 79	0

All (53) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	318	THR	4.0
2	D	267	VAL	3.9
2	C	355	TRP	3.7
2	C	359	LEU	3.5
2	D	302	PRO	3.4
2	C	358	VAL	3.3
2	C	302	PRO	3.2
2	D	359	LEU	3.2
2	C	269	LEU	3.2
2	C	267	VAL	3.2
2	C	301	GLY	3.1
2	C	300	ILE	2.9
1	B	128	LYS	2.9
2	D	355	TRP	2.8
2	D	301	GLY	2.8
2	C	327	HIS	2.8
2	C	299	LYS	2.8
2	D	299	LYS	2.8
2	C	357	THR	2.8
2	D	320	ASP	2.8
2	D	333	PHE	2.8

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Mol	Chain	Res	Type	RSRZ
2	C	295	VAL	2.7
2	D	358	VAL	2.7
2	C	268	ALA	2.7
2	C	320	ASP	2.7
2	D	263	ALA	2.7
2	D	303	ASP	2.6
2	D	295	VAL	2.6
2	C	356	LEU	2.5
2	C	333	PHE	2.5
2	C	329	ARG	2.5
2	D	269	LEU	2.5
1	B	46	VAL	2.5
2	D	328	LEU	2.5
2	D	300	ILE	2.4
2	D	330	ASN	2.4
2	D	327	HIS	2.3
2	D	325	VAL	2.3
2	D	357	THR	2.3
2	D	329	ARG	2.3
2	C	326	LEU	2.3
2	D	268	ALA	2.3
2	D	314	ALA	2.2
2	D	304	ASN	2.2
2	C	305	LEU	2.2
2	C	328	LEU	2.1
2	C	331	VAL	2.1
2	D	331	VAL	2.1
1	A	119	GLU	2.1
2	C	266	THR	2.1
2	C	304	ASN	2.1
2	C	298	SER	2.0
2	D	326	LEU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

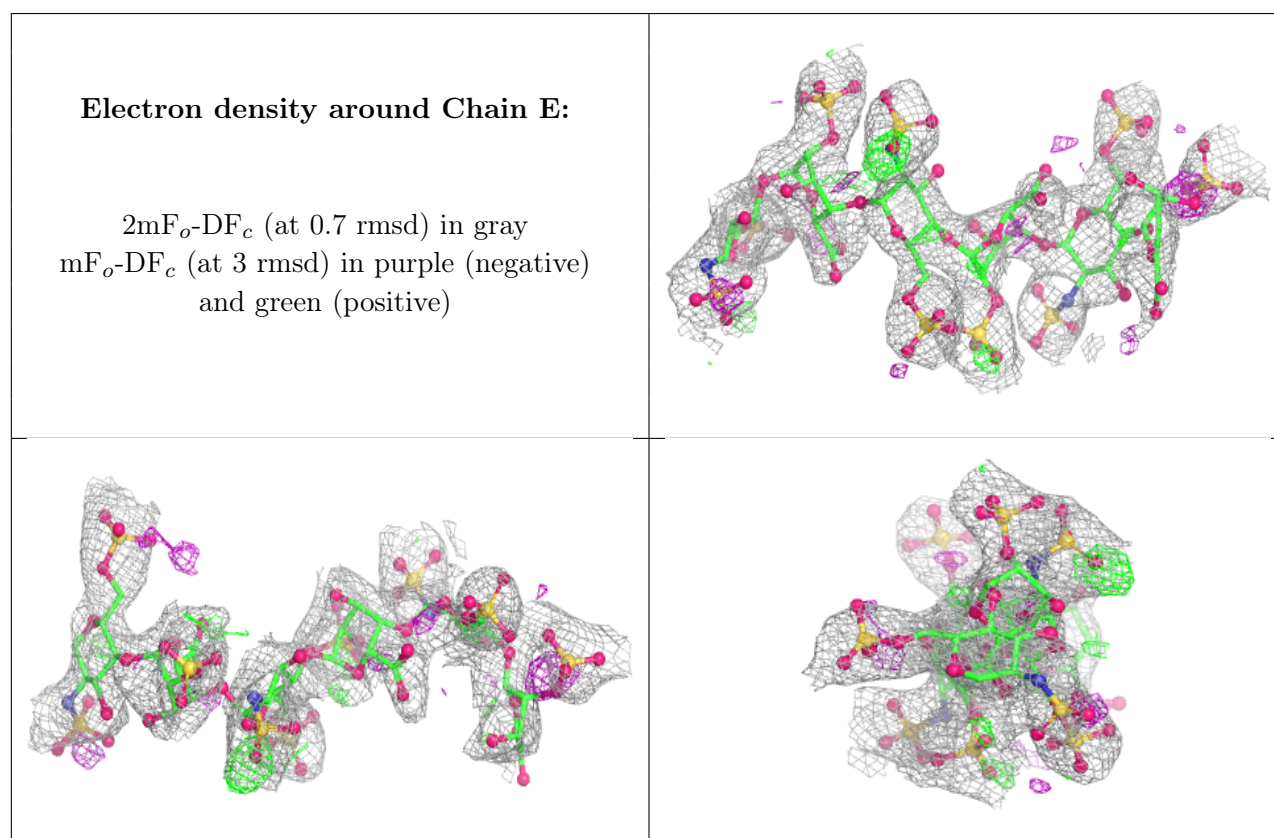
6.3 Carbohydrates [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum,

median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	UAP	E	6	15/16	0.48	0.20	84,86,87,87	0
3	SGN	E	1	19/20	0.60	0.19	87,90,94,94	0
3	SGN	E	5	19/20	0.75	0.19	78,81,85,85	0
3	IDS	E	4	16/17	0.75	0.17	74,79,79,80	0
3	SGN	E	3	19/20	0.76	0.16	78,79,81,81	0
3	IDS	E	2	16/17	0.77	0.19	82,83,84,85	0

The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.



6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.